

Data Path : Z:\HPCHEM1\BNA G\DATA\BG030316\
 Data File : BG021251.D
 Acq On : 4 Mar 2016 2:12
 Operator : UM/SJ
 Sample : H1626-11
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 JHFJ8

Quant Time: Mar 04 07:17:18 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG030316.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 04 02:35:52 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.13	152	10327	20.00	ng/ul	0.00
18) Naphthalene-d8	10.94	136	50603	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.74	164	28712	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.48	188	62793	20.00	ng/ul	0.00
78) Chrysene-d12	21.74	240	71780	20.00	ng/ul	0.00
86) Perylene-d12	24.99	264	66010	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.55	96	386	1.52	ng/uL	0.00
5) Phenol-d5	7.28	99	9280	8.48	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.45	67	25209	34.56	ng/ul	0.00
9) 2-Chlorophenol-d4	7.66	132	22233	29.13	ng/ul	0.00
13) 4-Methylphenol-d8	8.82	113	15806	18.30	ng/ul	0.00
19) Nitrobenzene-d5	9.29	128	14120	34.45	ng/ul	0.00
22) 2-Nitrophenol-d4	10.01	143	14900	34.10	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.55	165	22862	29.28	ng/ul	0.00
29) 4-Chloroaniline-d4	11.08	131	1529	1.54	ng/ul	0.00
44) Dimethylphthalate-d6	14.14	166	82683	34.36	ng/ul	0.00
47) Acenaphthylene-d8	14.43	160	105653	35.29	ng/ul	0.00
52) 4-Nitrophenol-d4	14.91	143	3325	7.09	ng/ul	0.00
58) Fluorene-d10	15.73	176	73086	34.59	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.83	200	13904	30.85	ng/ul	0.00
71) Anthracene-d10	17.58	188	111252	35.63	ng/ul	0.00
79) Pyrene-d10	19.85	212	128893	35.47	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.77	264	125582	35.24	ng/ul	0.00

Target Compounds Ovalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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